

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1617srh

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page for STN Seminar Schedule - N. America
NEWS 2 AUG 15 CAOLD to be discontinued on December 31, 2008
NEWS 3 OCT 07 EPFULL enhanced with full implementation of EPC2000
NEWS 4 OCT 07 Multiple databases enhanced for more flexible patent
number searching
NEWS 5 OCT 22 Current-awareness alert (SDI) setup and editing
enhanced
NEWS 6 OCT 22 WPIDS, WPINDEX, and WPIX enhanced with Canadian PCT
Applications
NEWS 7 OCT 24 CHEMLIST enhanced with intermediate list of
pre-registered REACH substances
NEWS 8 NOV 21 CAS patent coverage to include exemplified prophetic
substances identified in English-, French-, German-,
and Japanese-language basic patents from 2004-present
NEWS 9 NOV 26 MARPAT enhanced with FSORT command
NEWS 10 NOV 26 MEDLINE year-end processing temporarily halts
availability of new fully-indexed citations
NEWS 11 NOV 26 CHEMSAFE now available on STN Easy
NEWS 12 NOV 26 Two new SET commands increase convenience of STN
searching
NEWS 13 DEC 01 ChemPort single article sales feature unavailable
NEWS 14 DEC 12 GBFULL now offers single source for full-text
coverage of complete UK patent families
NEWS 15 DEC 17 Fifty-one pharmaceutical ingredients added to PS

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS LOGIN Welcome Banner and News Items
NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that
specific topic.

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of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 15:24:00 ON 05 JAN 2009

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

2.20

2.20

FILE 'REGISTRY' ENTERED AT 15:29:59 ON 05 JAN 2009

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 4 JAN 2009 HIGHEST RN 1092523-63-1

DICTIONARY FILE UPDATES: 4 JAN 2009 HIGHEST RN 1092523-63-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

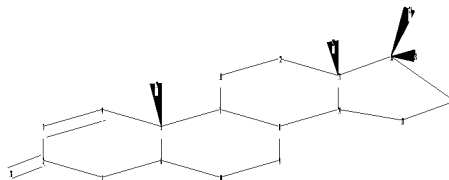
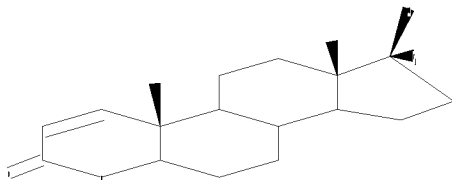
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10-501640 brd.str



chain nodes :
18 19 20 21 23
ring nodes :

```

1  2  3  4  5  6  7  8  9  10  11  12  13  14  15  16  17
chain bonds :
2-18  5-19  13-20  15-21  15-23
ring bonds :
1-2  1-6  2-3  3-4  4-5  5-6  5-7  6-10  7-8  7-11  8-9  8-14  9-10  11-12  12-13
13-14  13-15  14-17  15-16  16-17
exact/norm bonds :
1-2  1-6  2-3  2-18  3-4  4-5  5-6  5-7  6-10  7-8  7-11  8-9  8-14  9-10  11-12
12-13  13-14  13-15  14-17  15-16  15-21  15-23  16-17
exact bonds :
5-19  13-20

```

G1:H,Ak

```

Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom  9:Atom  10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:CLASS 19:CLASS
20:CLASS 21:CLASS 23:CLASS

```

Stereo Bonds:

```

19-5 (Single Wedge).
20-13 (Single Wedge).
21-15 (Single Wedge).
23-15 (Single Hash).

```

Stereo Chiral Centers:

```

5      (Parity=Even)
13     (Parity=Even)
15     (Parity=Odd)

```

Stereo RSS Sets:

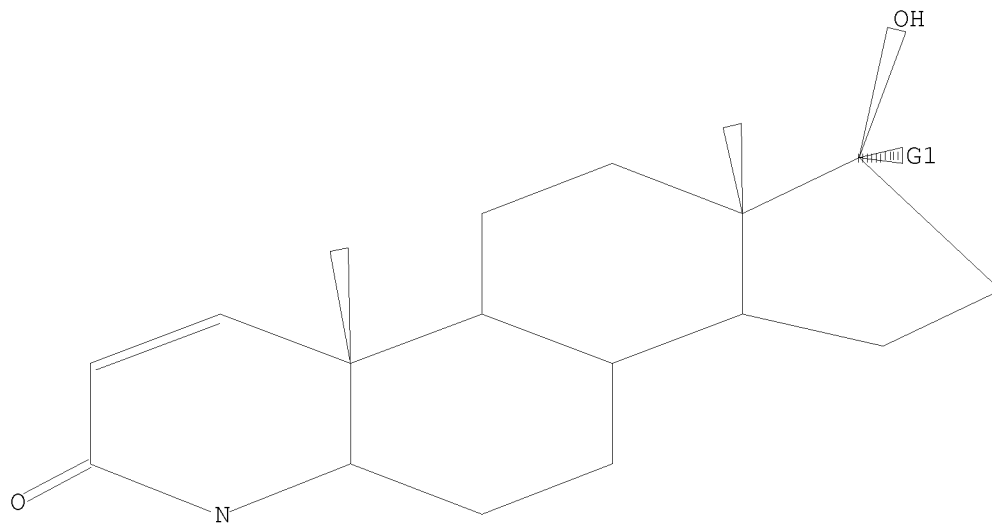
Type=Relative (Default). 3 Nodes= 5 13 15

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR



G1 H,Ak

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 15:30:21 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 411 TO ITERATE

100.0% PROCESSED 411 ITERATIONS

3 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 7004 TO 9436

PROJECTED ANSWERS: 3 TO 163

L2 3 SEA SSS SAM L1

=> d scan

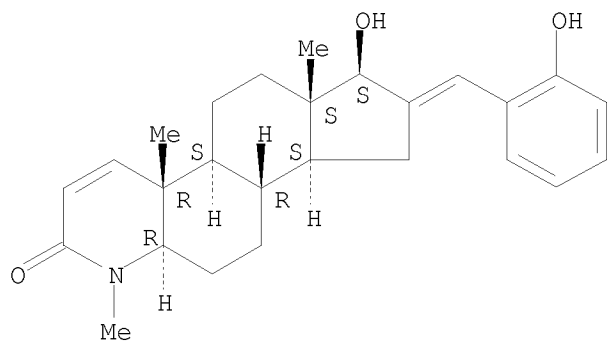
L2 3 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-8-[(2-hydroxyphenyl)methylene]-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-

MF C26 H33 N O3

Absolute stereochemistry.

Double bond geometry unknown.

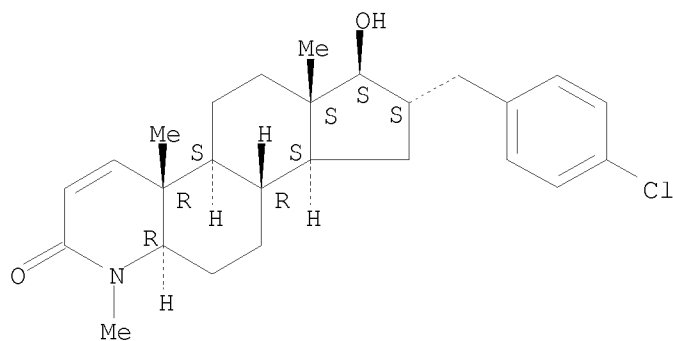


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L2 3 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(4-chlorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H34 Cl N O2

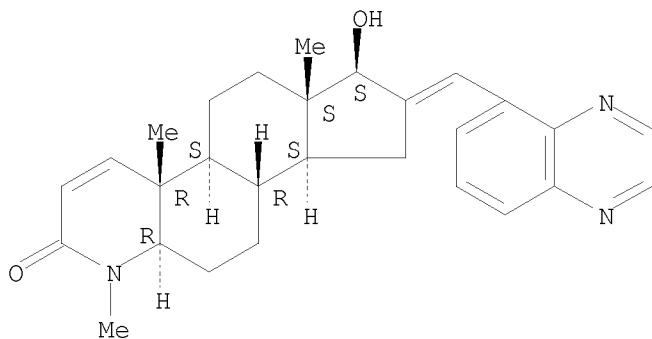
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 3 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(5-quinoxalinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H33 N3 O2

Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> s l1 full
 FULL SEARCH INITIATED 15:31:52 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 7600 TO ITERATE

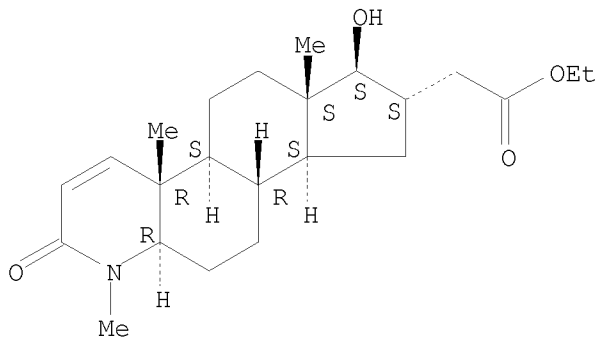
100.0% PROCESSED 7600 ITERATIONS 103 ANSWERS
 SEARCH TIME: 00.00.01

L3 103 SEA SSS FUL L1

=> d scan

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 1H-Indeno[5,4-f]quinoline-8-acetic acid,
 2,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-2-oxo-, ethyl ester, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C23 H35 N O4

Absolute stereochemistry.

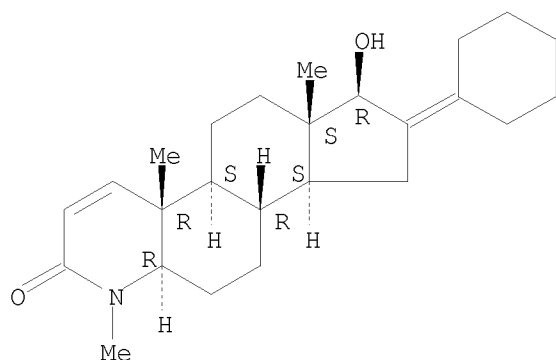


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):102

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-cyclohexylidene-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 MF C25 H37 N O2

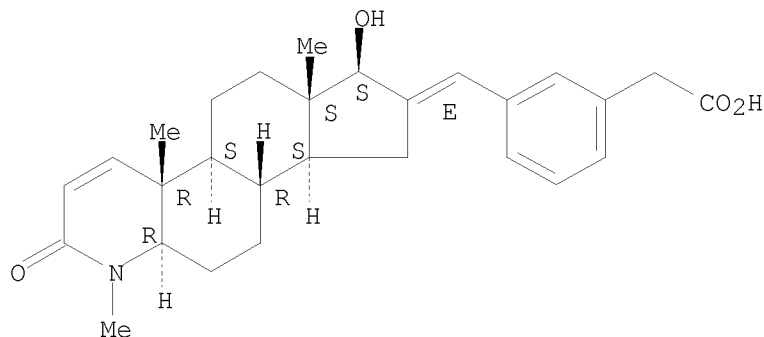
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN Benzeneacetic acid, 3-[(E)-[(4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 1,2,4a,4b,5,6,6a,7,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-2-oxo-8H-indeno[5,4-f]quinolin-8-ylidene)methyl]-
 MF C28 H35 N O4

Absolute stereochemistry.
 Double bond geometry as shown.

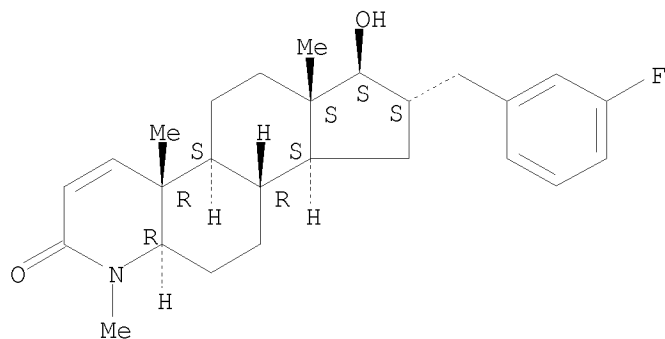


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3-fluorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-

trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
MF C26 H34 F N O2

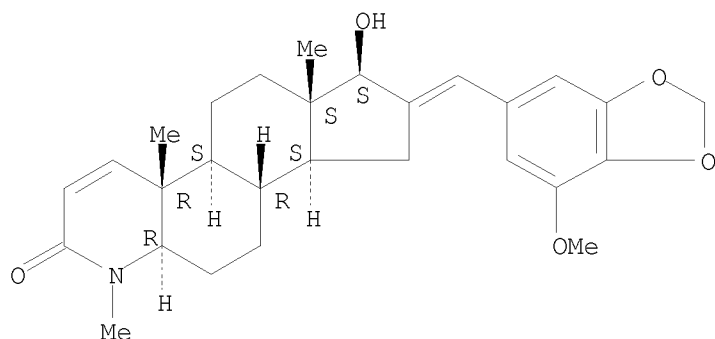
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
tetradecahydro-7-hydroxy-8-[(7-methoxy-1,3-benzodioxol-5-yl)methylene]-
1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
MF C28 H35 N O5

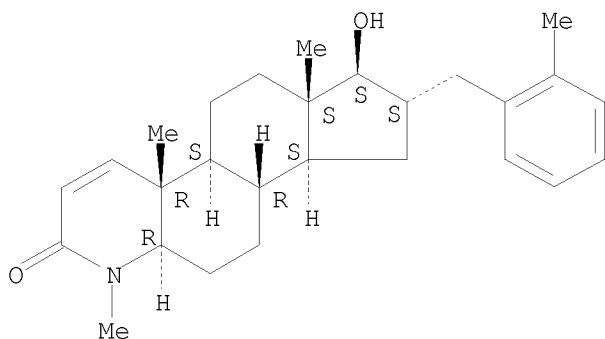
Absolute stereochemistry.
Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(2-methylphenyl)methyl]-,
(4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
MF C27 H37 N O2

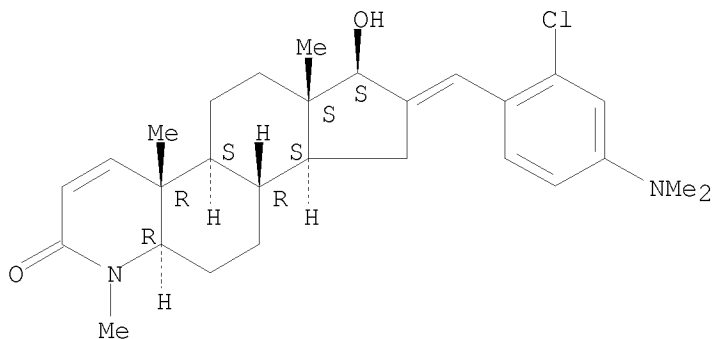
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[[2-chloro-4-(dimethylamino)phenyl]methylene]-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H37 Cl N2 O2

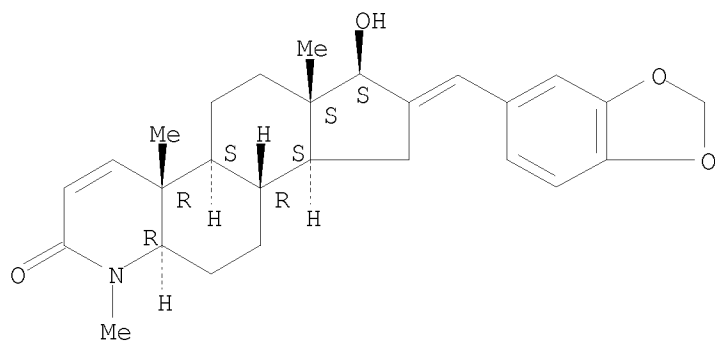
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-(1,3-benzodioxol-5-ylmethylene)-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H33 N O4

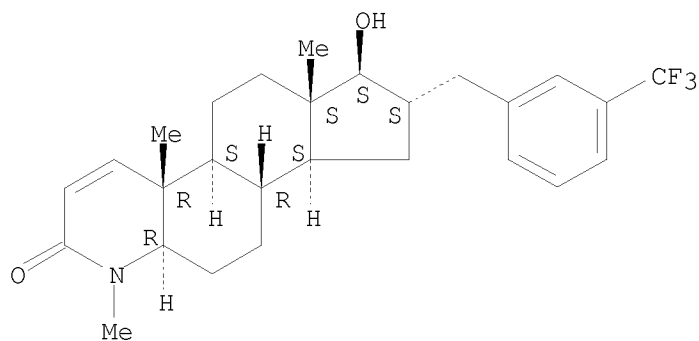
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[3-
 (trifluoromethyl)phenyl]methyl]-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C27 H34 F3 N O2

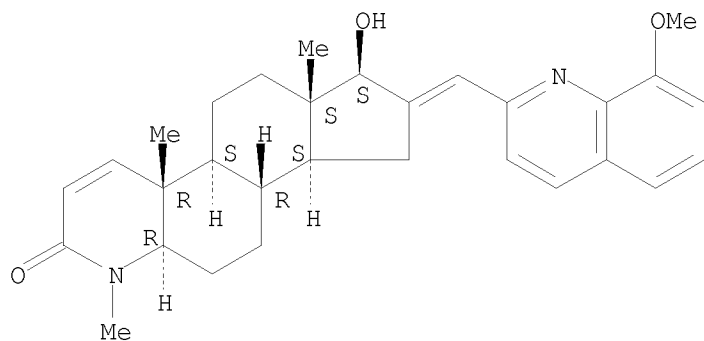
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(8-methoxy-2-quinolinyl)methylene]-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C30 H36 N2 O3

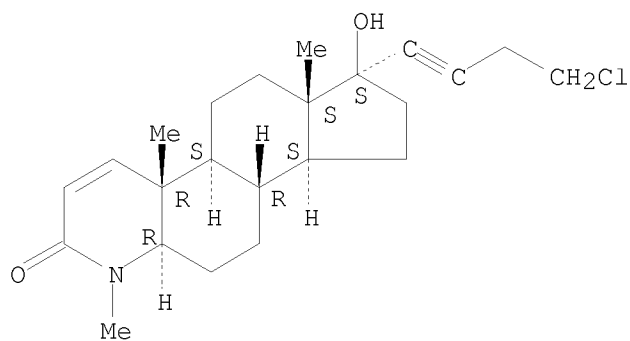
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 7-(4-chloro-1-butyne-1-yl)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H32 Cl N O2

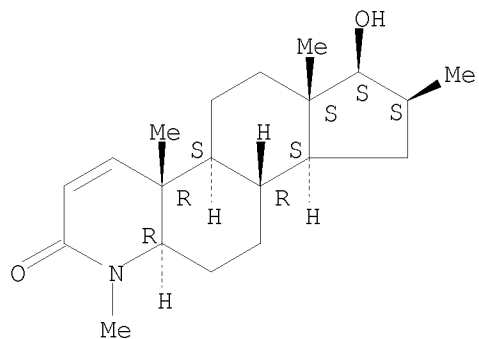
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a,8-tetramethyl-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C20 H31 N O2

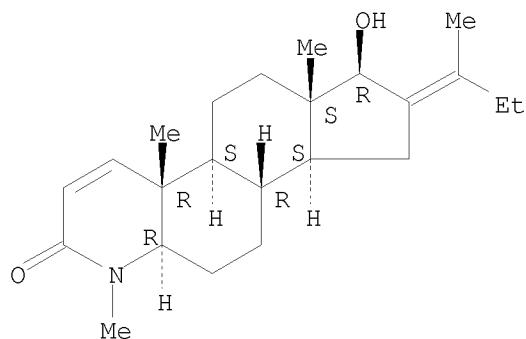
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(1-methylpropylidene)-,
 (4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 MF C23 H35 N O2

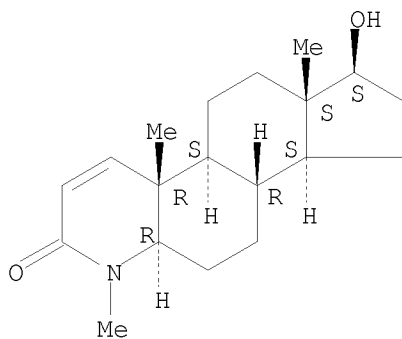
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C19 H29 N O2

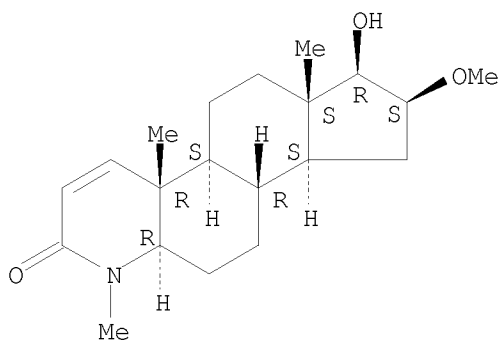
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-methoxy-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7R,8S,9aS,9bR,11aR)-
 MF C20 H31 N O3

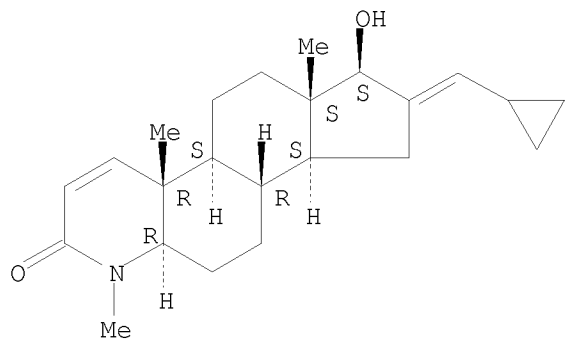
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-(cyclopropylmethylene)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H33 N O2

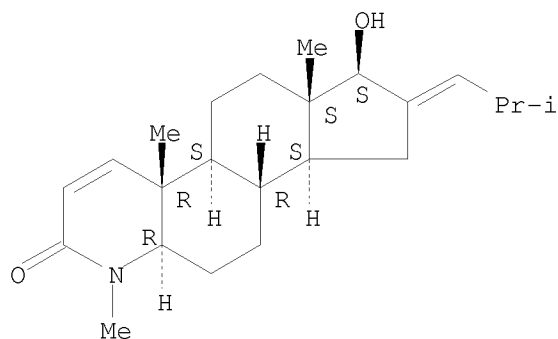
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(2-methylpropylidene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H35 N O2

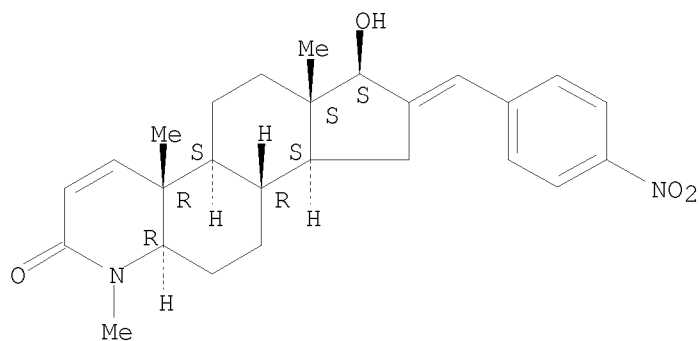
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(4-nitrophenyl)methylene]-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H32 N2 O4

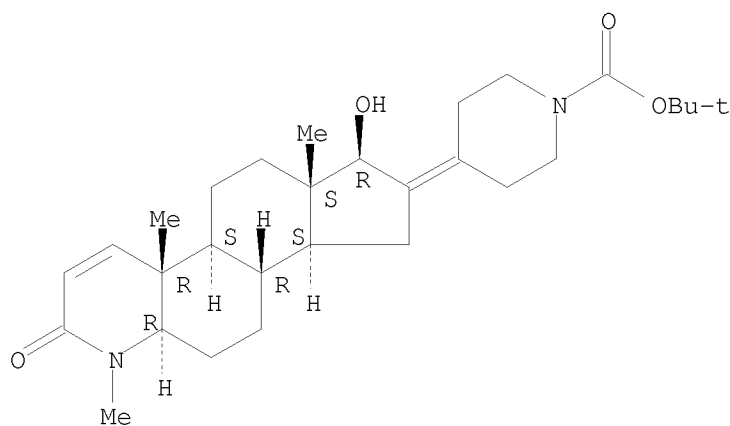
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 1-Piperidinecarboxylic acid, 4-[(4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 1,2,4a,4b,5,6,6a,7,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-2-oxo-8H-indeno[5,4-f]quinolin-8-ylidene]-, 1,1-dimethylethyl
 ester
 MF C29 H44 N2 O4

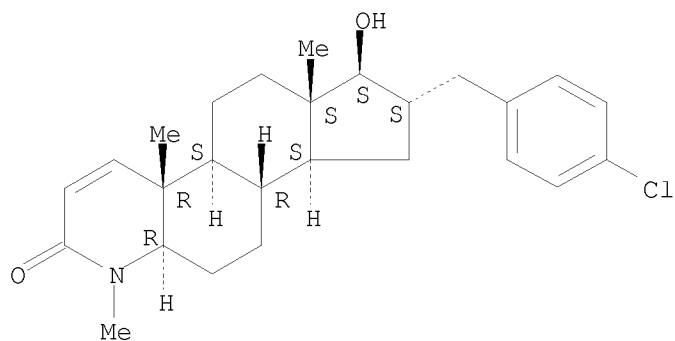
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(4-chlorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H34 Cl N O2

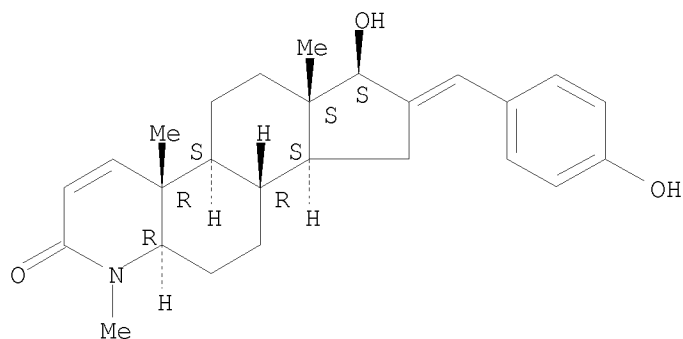
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(4-hydroxyphenyl)methylene]-1,4a,6a-trimethyl-
 , (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H33 N O3

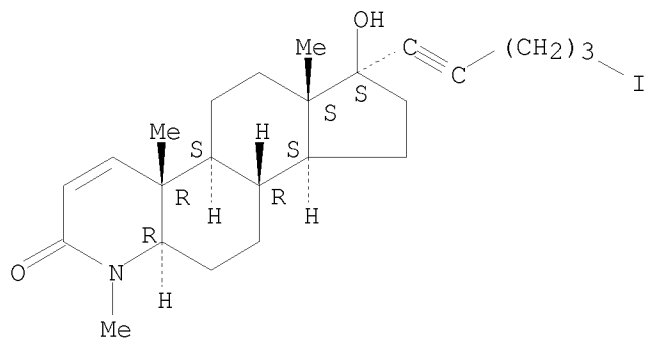
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-7-(5-iodo-1-pentyn-1-yl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H34 I N O2

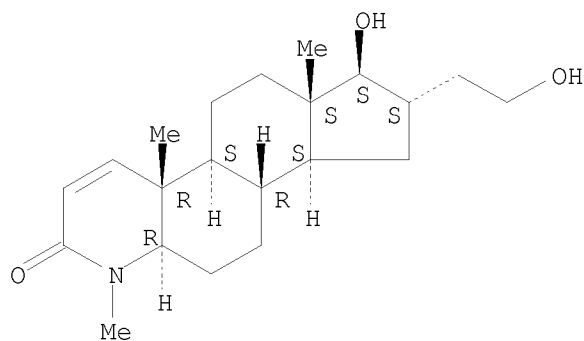
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-(2-hydroxyethyl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C21 H33 N O3

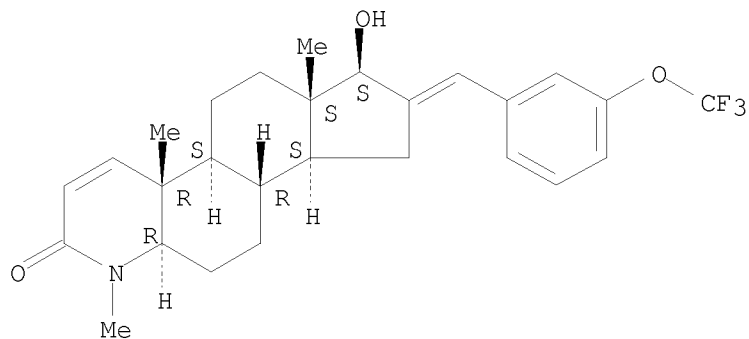
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[3-
 (trifluoromethoxy)phenyl]methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H32 F3 N O3

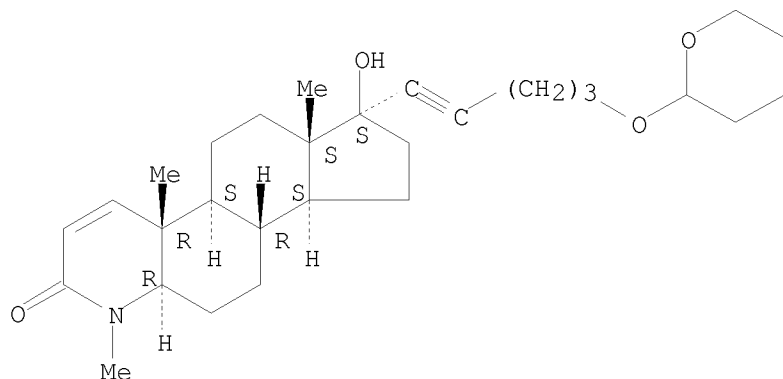
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-7-[5-[(tetrahydro-2H-pyran-2-yl)oxy]-1-pentyn-1-yl]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C29 H43 N O4

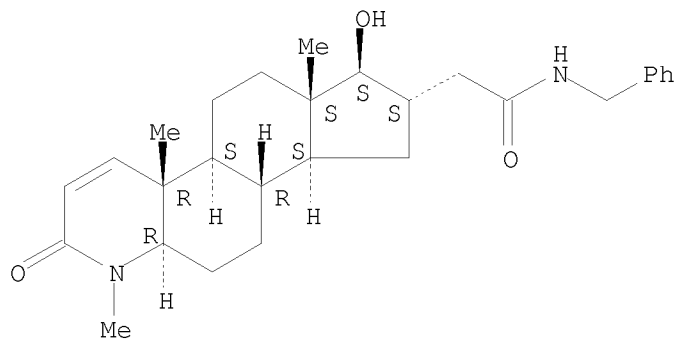
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 1H-Indeno[5,4-f]quinoline-8-acetamide, 2,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-2-oxo-N-(phenylmethyl)-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C28 H38 N2 O3

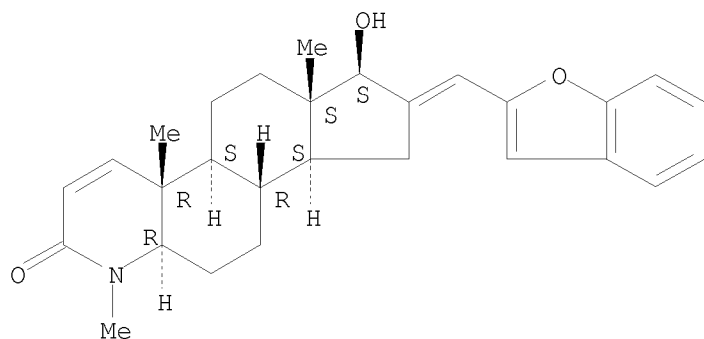
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-(2-benzofuranylmethylene)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H33 N O3

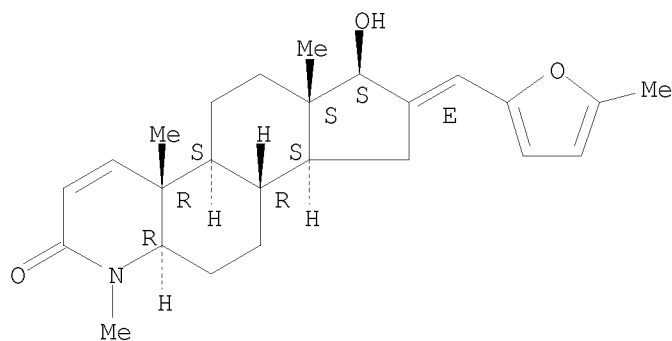
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(5-methyl-2-
 furanyl)methylene]-, (4aR,4bS,6aS,7S,8E,9aS,9bR,11aR)-
 MF C25 H33 N O3

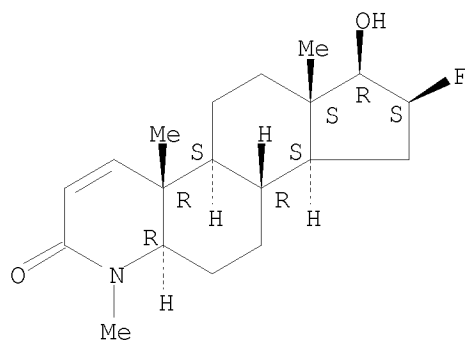
Absolute stereochemistry.
 Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-fluoro-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7R,8S,9aS,9bR,11aR)-
 MF C19 H28 F N O2

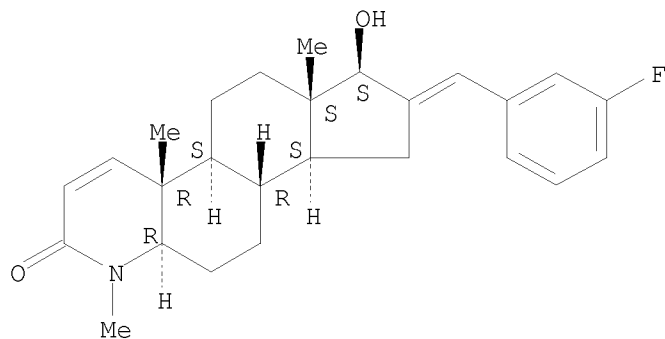
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3-fluorophenyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H32 F N O2

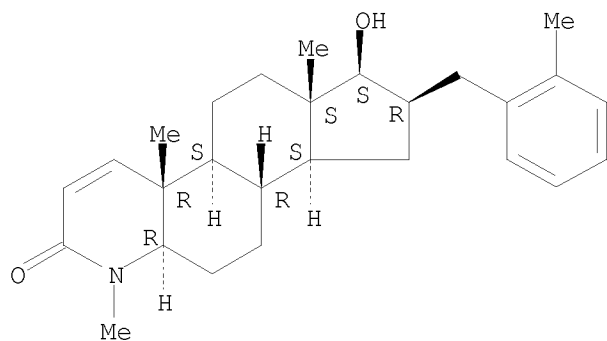
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(2-methylphenyl)methyl]-, (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C27 H37 N O2

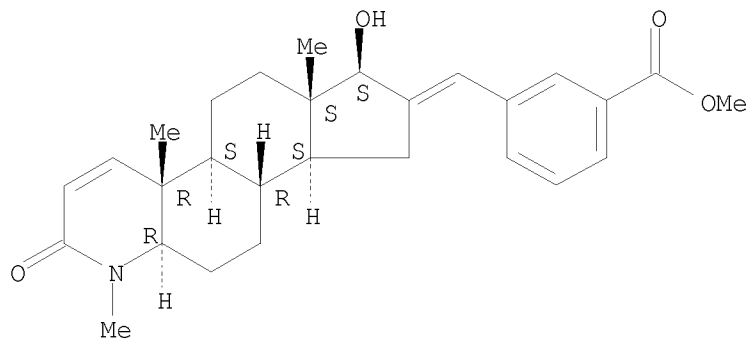
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN Benzoic acid, 3-[[[(4aR,4bS,6aS,7S,9aS,9bR,11aR)-1,2,4a,4b,5,6,6a,7,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-2-oxo-8H-indeno[5,4-f]quinolin-8-ylidene)methyl]-, methyl ester
 MF C28 H35 N O4

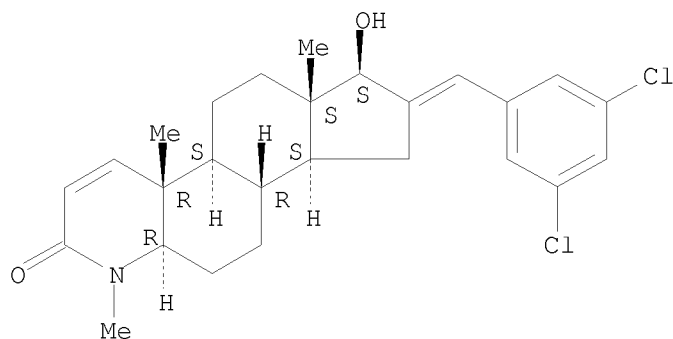
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3,5-dichlorophenyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H31 Cl2 N O2

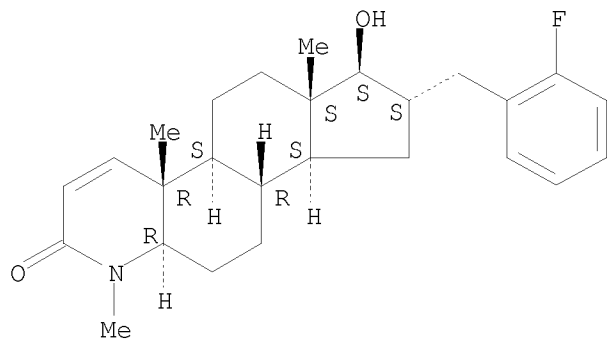
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(2-fluorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H34 F N O2

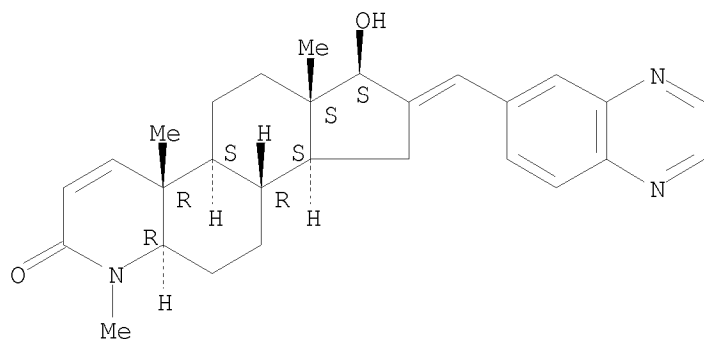
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(6-quinoxalinylmethylene)-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H33 N3 O2

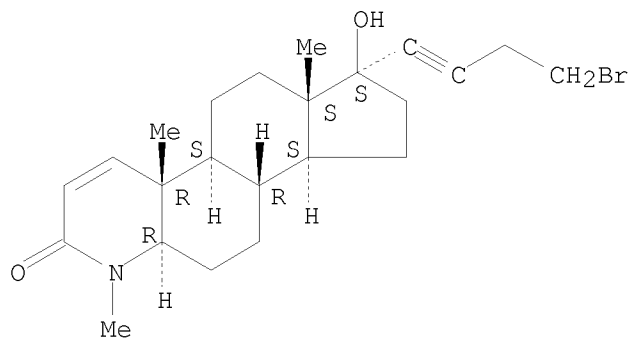
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 7-(4-bromo-1-butyn-1-yl)-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H32 Br N O2

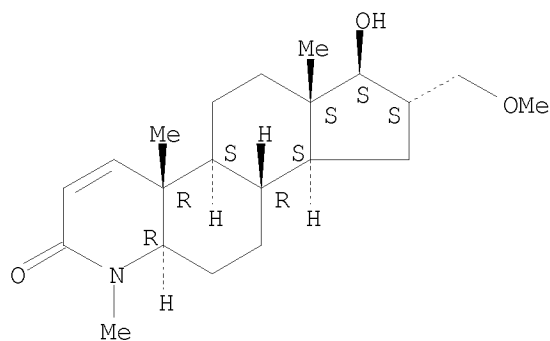
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-(methoxymethyl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C21 H33 N O3

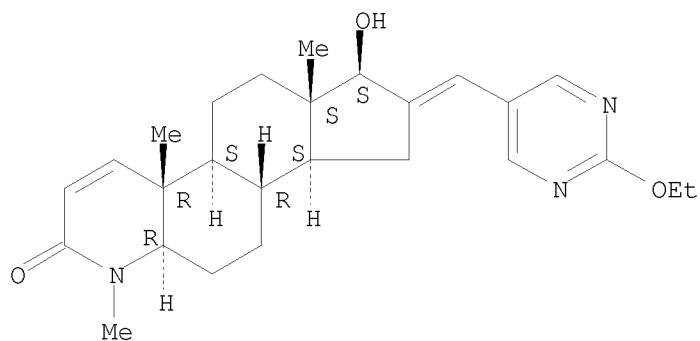
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(2-ethoxy-5-pyrimidinyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H35 N3 O3

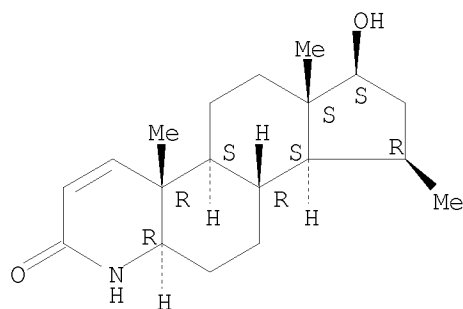
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 2,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-4a,6a,9-trimethyl-,
 (4aR,4bS,6aS,7S,9R,9aS,9bR,11aR)-(9CI)
 MF C19 H29 N O2

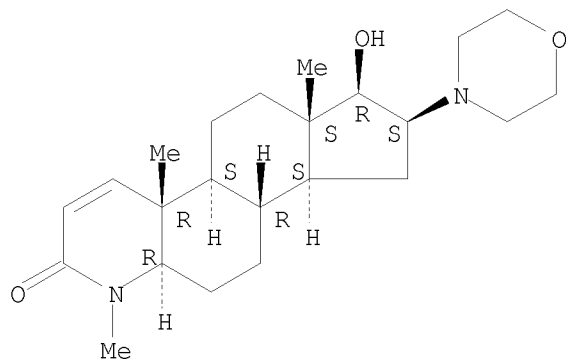
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(4-morpholinyl)-,
 (4aR,4bS,6aS,7R,8S,9aS,9bR,11aR)-
 MF C23 H36 N2 O3

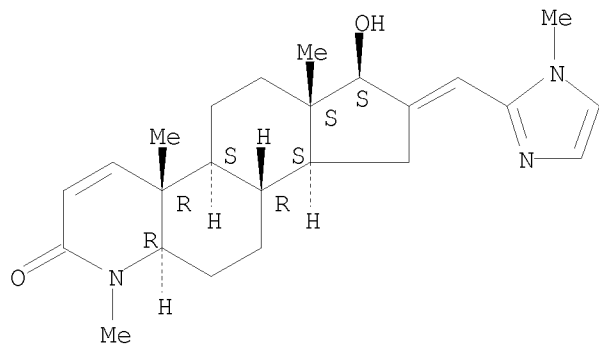
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(1-methyl-1H-imidazol-2-yl)methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H33 N3 O2

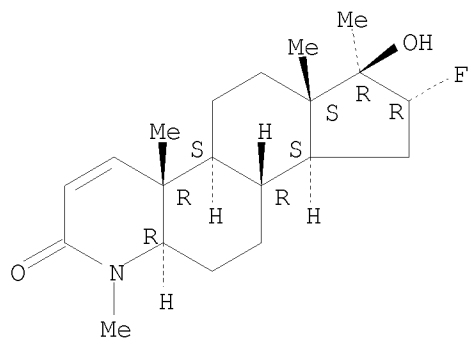
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-fluoro-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a,7-tetramethyl-, (4aR,4bS,6aS,7R,8R,9aS,9bR,11aR)-
 MF C20 H30 F N O2

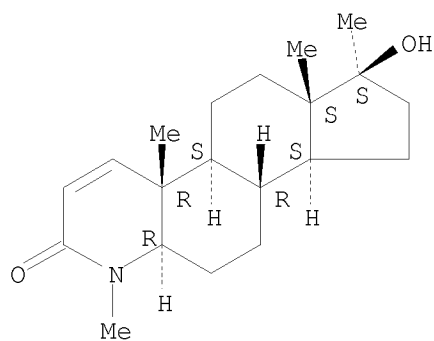
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a,7-tetramethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C20 H31 N O2

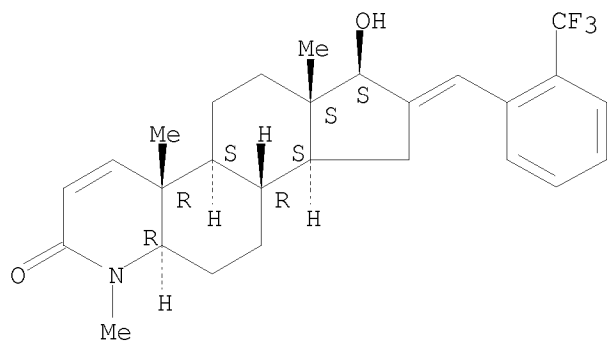
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[2-
 (trifluoromethyl)phenyl]methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H32 F3 N O2

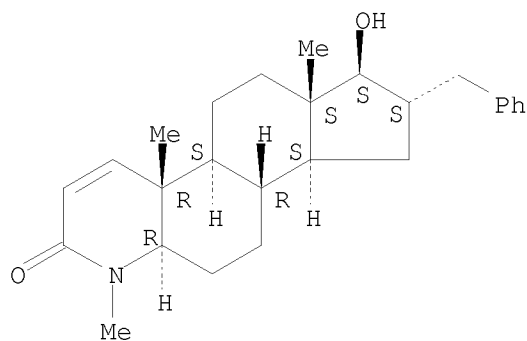
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(phenylmethyl)-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H35 N O2

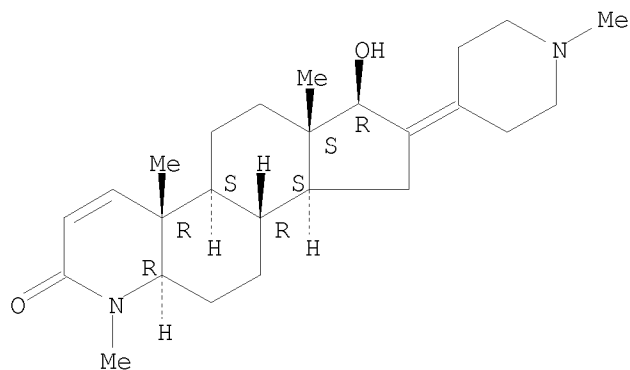
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(1-methyl-4-piperidinyldene)-
 , (4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 MF C25 H38 N2 O2

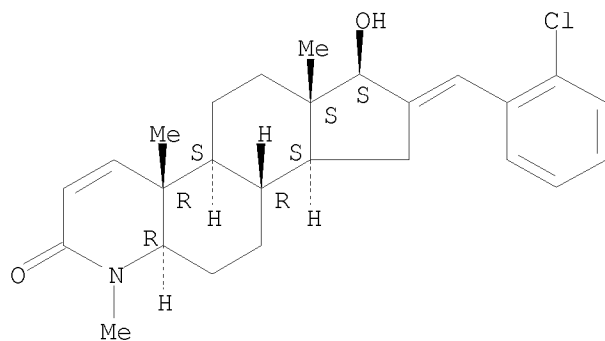
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(2-chlorophenyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H32 Cl N O2

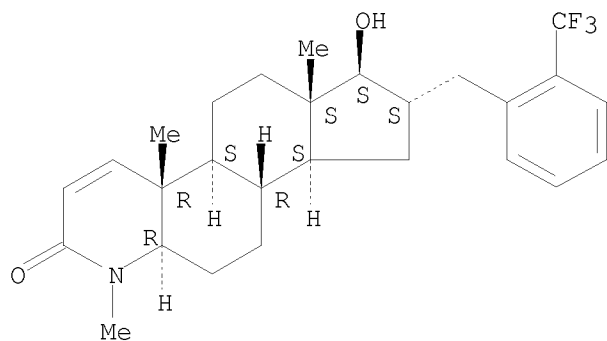
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[2-
 (trifluoromethyl)phenyl]methyl]-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C27 H34 F3 N O2

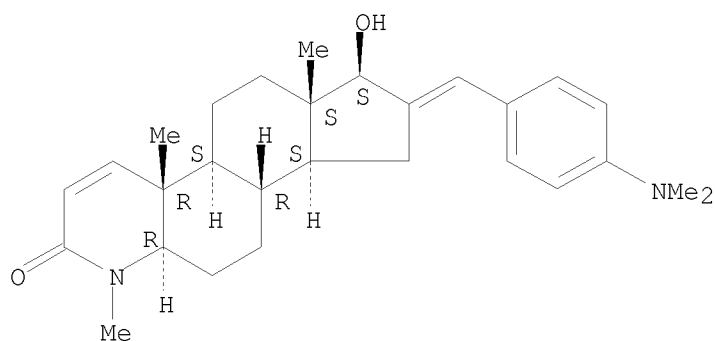
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[[4-(dimethylamino)phenyl]methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H38 N2 O2

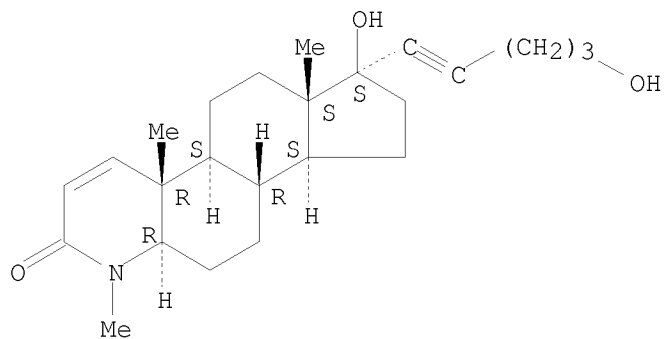
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-7-(5-hydroxy-1-pentyn-1-yl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H35 N O3

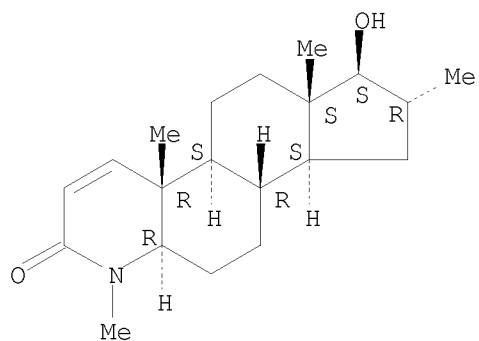
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a,8-tetramethyl-,
 (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C20 H31 N O2

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(1-methyl-1H-pyrrol-2-
 yl)methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H34 N2 O2

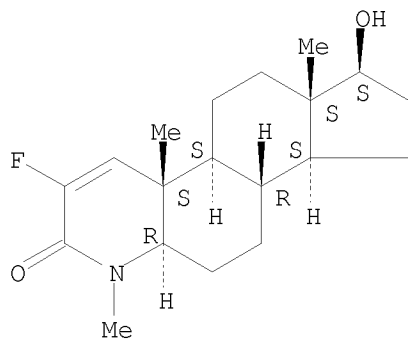
Absolute stereochemistry.
 Double bond geometry unknown.



Absolute stereochemistry.



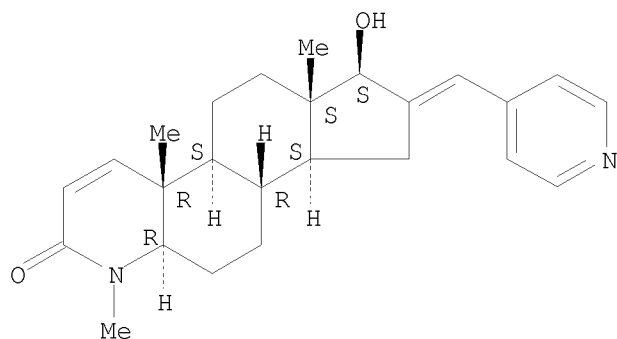
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(4-pyridinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H32 N2 O2

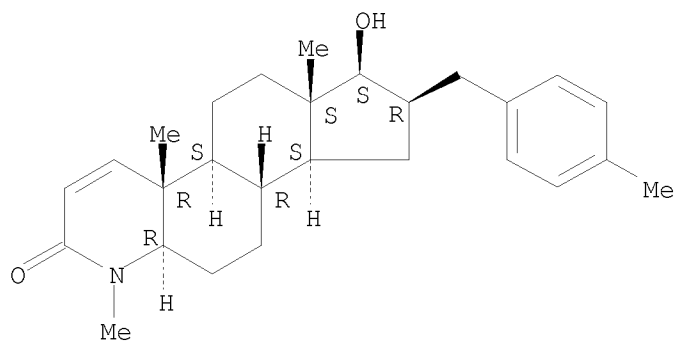
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(4-methylphenyl)methyl]-,
 (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C27 H37 N O2

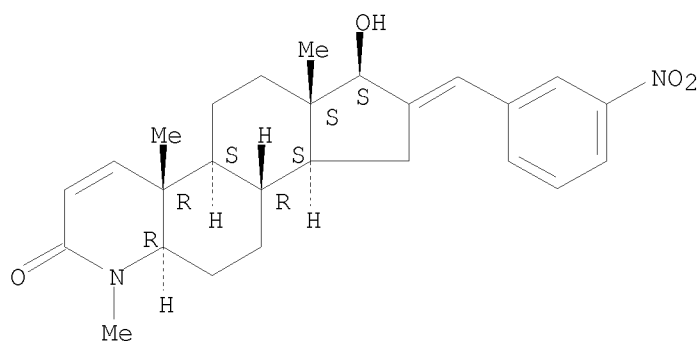
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(3-nitrophenyl)methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H32 N2 O4

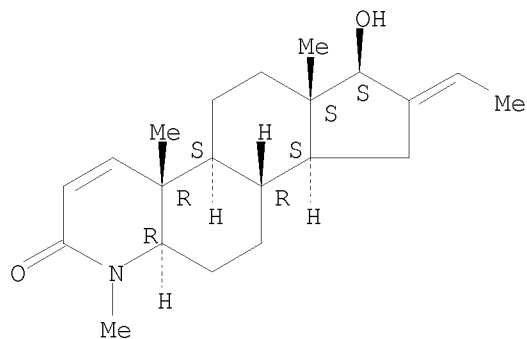
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-ethylidene-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C21 H31 N O2

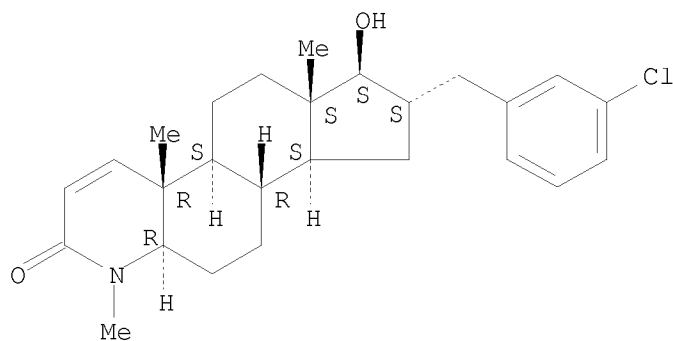
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3-chlorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H34 Cl N O2

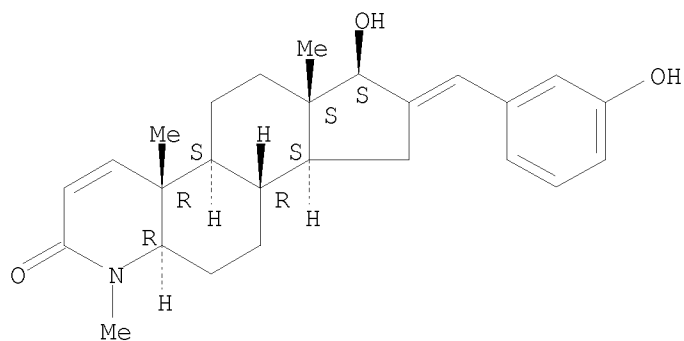
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(3-hydroxyphenyl)methylene]-1,4a,6a-trimethyl-
 , (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H33 N O3

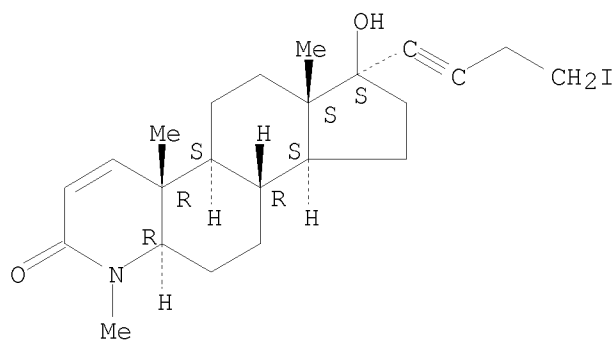
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-7-(4-iodo-1-butyn-1-yl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H32 I N O2

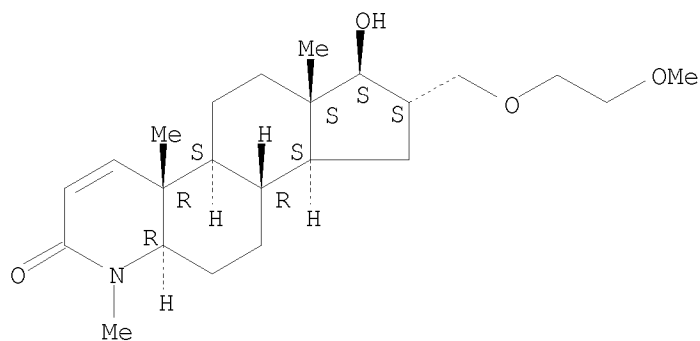
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(2-methoxyethoxy)methyl]-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C23 H37 N O4

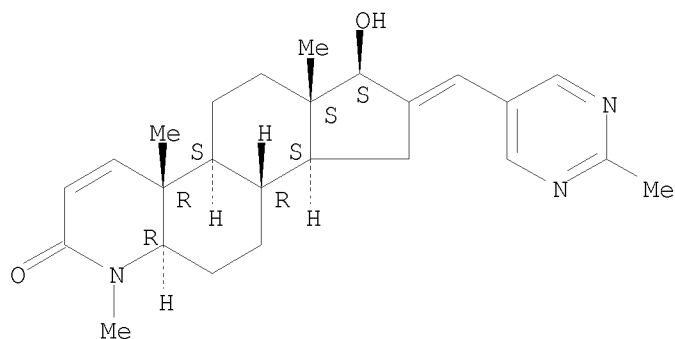
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(2-methyl-5-
 pyrimidinyl)methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H33 N3 O2

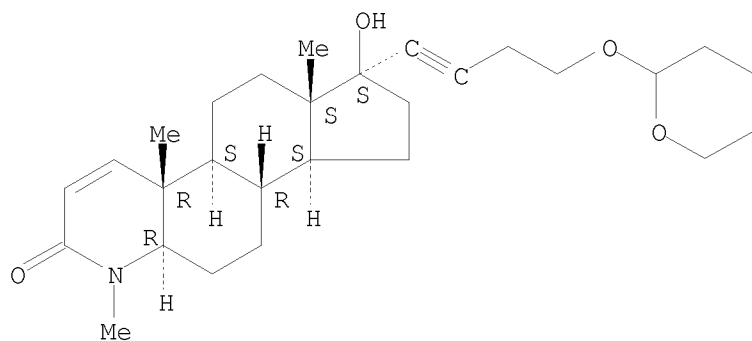
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-7-[4-[(tetrahydro-2H-pyran-2-
 yl)oxy]-1-butyn-1-yl]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H41 N O4

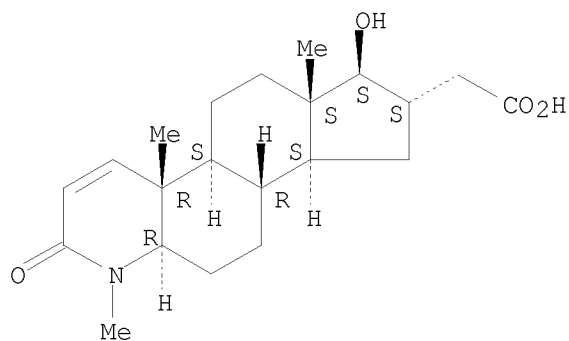
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 1H-Indeno[5,4-f]quinoline-8-acetic acid,
 2,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-2-oxo-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C21 H31 N O4

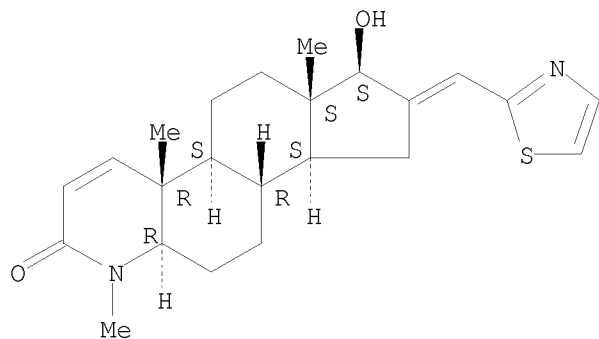
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(2-thiazolylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H30 N2 O2 S

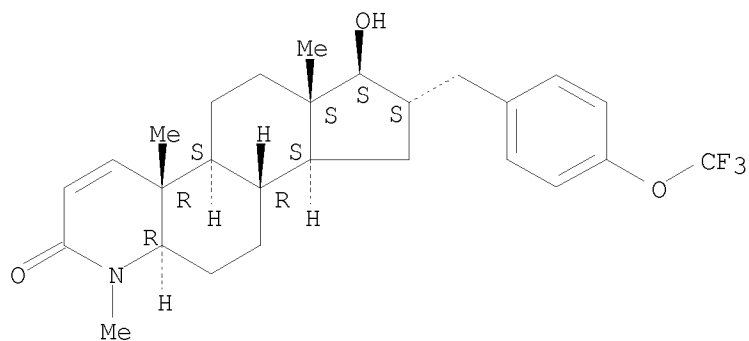
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[4-
 (trifluoromethoxy)phenyl]methyl]-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C27 H34 F3 N O3

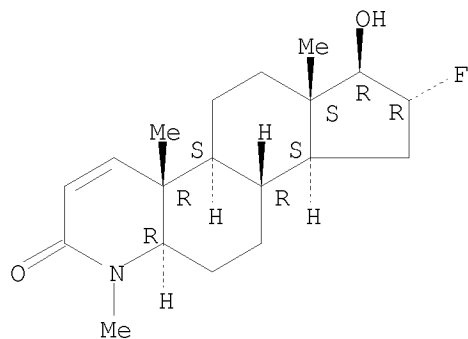
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-fluoro-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7R,8R,9aS,9bR,11aR)-
 MF C19 H28 F N O2

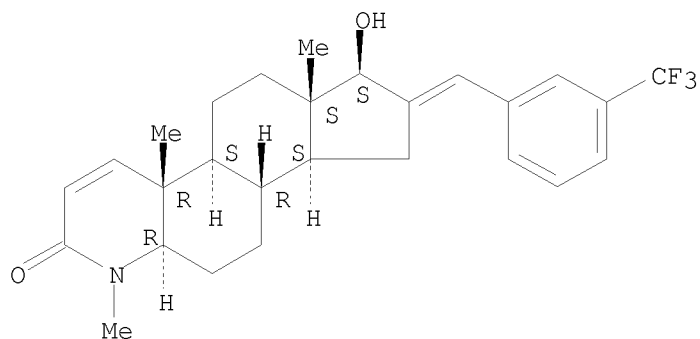
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[3-(trifluoromethyl)phenyl]methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H32 F3 N O2

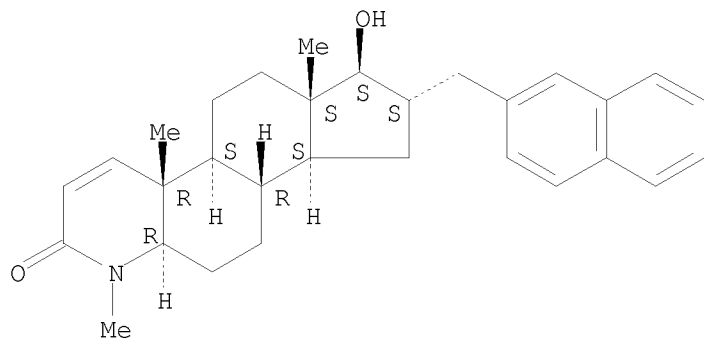
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(2-naphthalenylmethyl)-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C30 H37 N O2

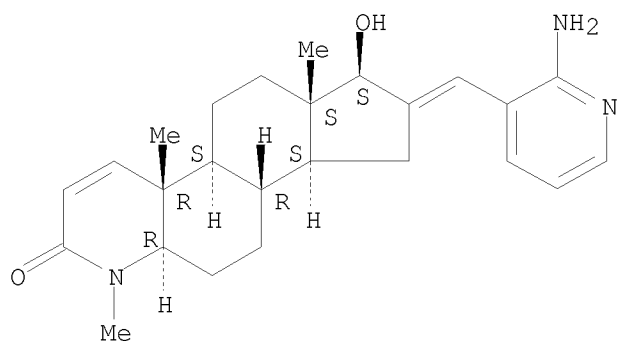
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(2-amino-3-pyridinyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H33 N3 O2

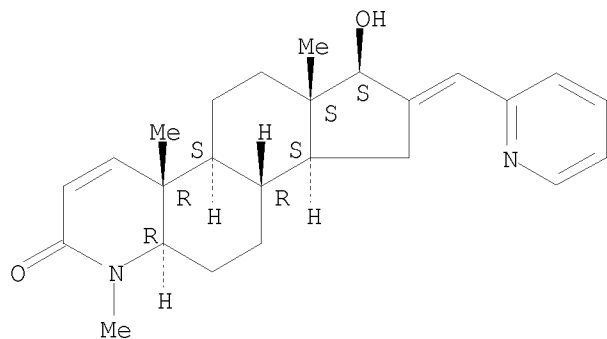
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(2-pyridinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H32 N2 O2

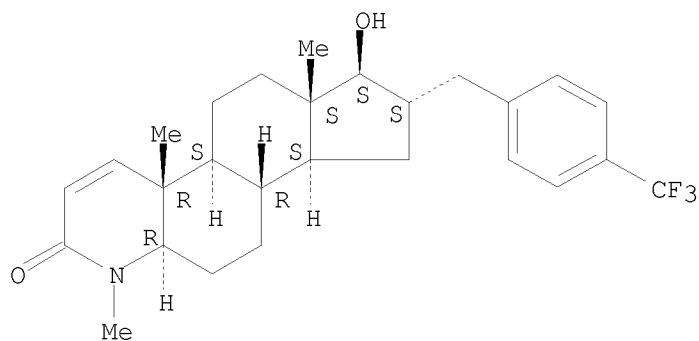
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[4-
 (trifluoromethyl)phenyl]methyl]-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C27 H34 F3 N O2

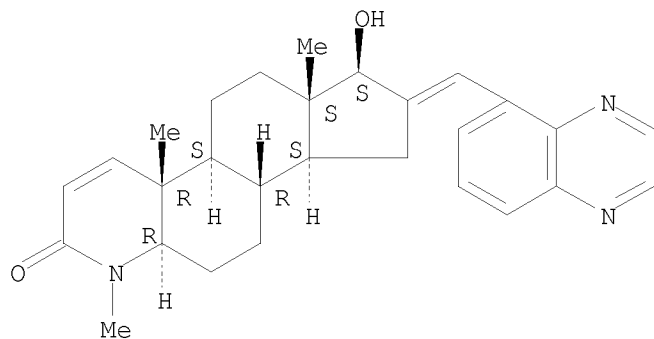
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(5-quinoxalinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H33 N3 O2

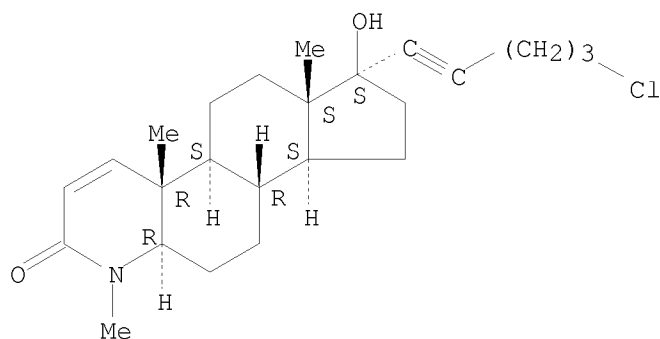
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 7-(5-chloro-1-pentyn-1-yl)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H34 Cl N O2

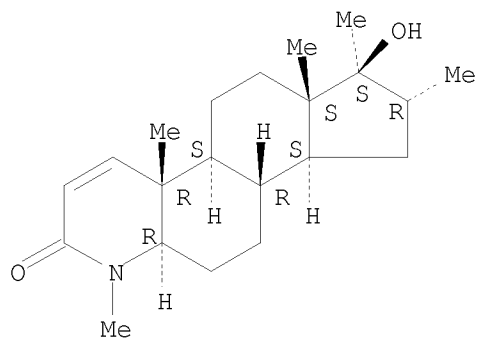
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a,7,8-pentamethyl-,
 (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C21 H33 N O2

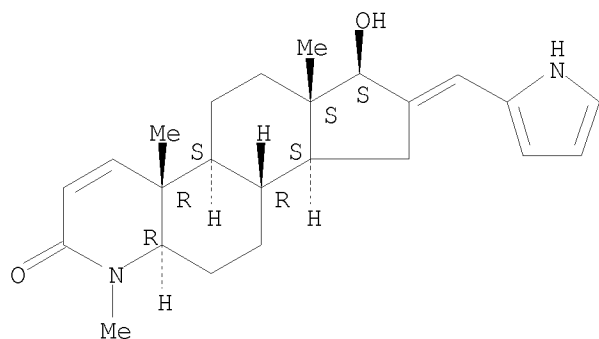
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(1H-pyrrol-2-ylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H32 N2 O2

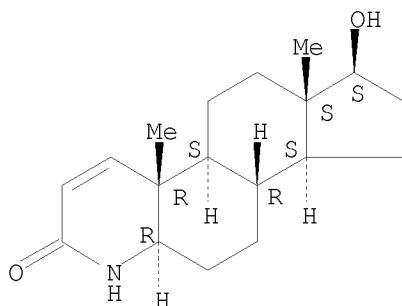
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-4a,6a-dimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C18 H27 N O2

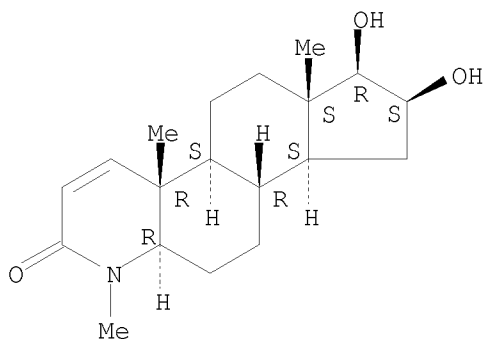
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7,8-dihydroxy-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7R,8S,9aS,9bR,11aR)-
 MF C19 H29 N O3

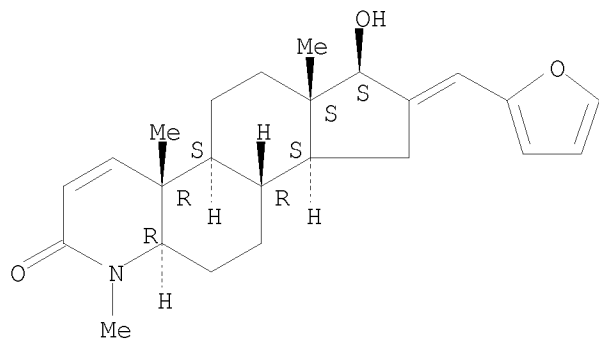
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-(2-furanylmethylene)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H31 N O3

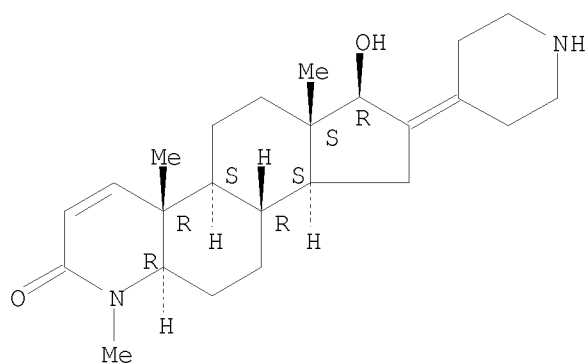
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(4-piperidinylidene)-,
 (4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 MF C24 H36 N2 O2

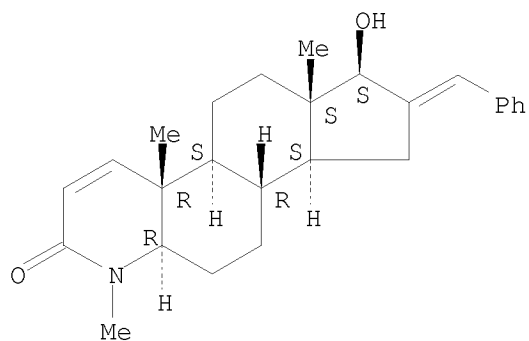
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(phenylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H33 N O2

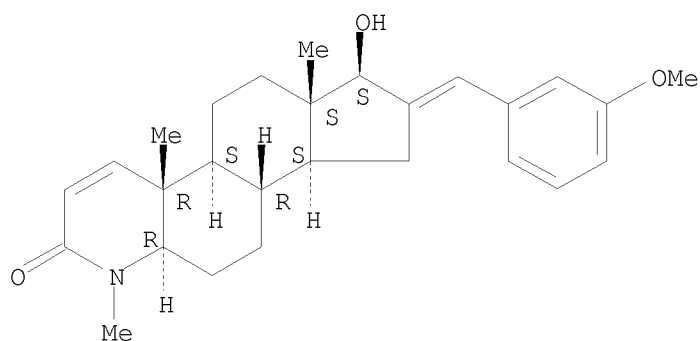
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(3-methoxyphenyl)methylene]-1,4a,6a-trimethyl-
 , (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H35 N O3

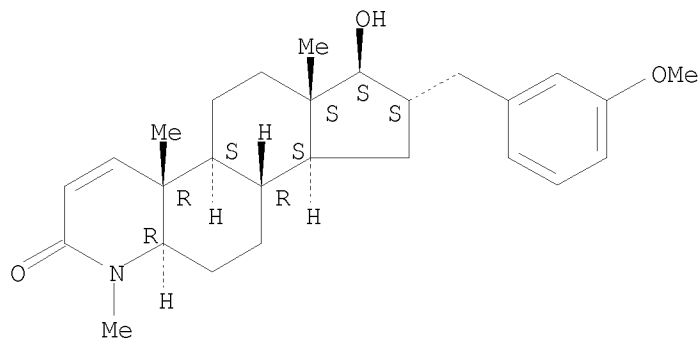
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[(3-methoxyphenyl)methyl]-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C27 H37 N O3

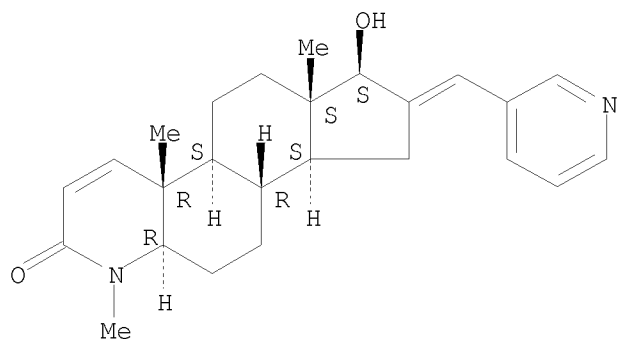
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(3-pyridinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H32 N2 O2

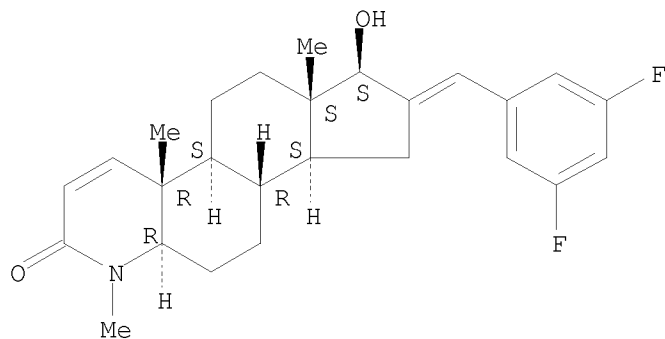
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3,5-difluorophenyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H31 F2 N O2

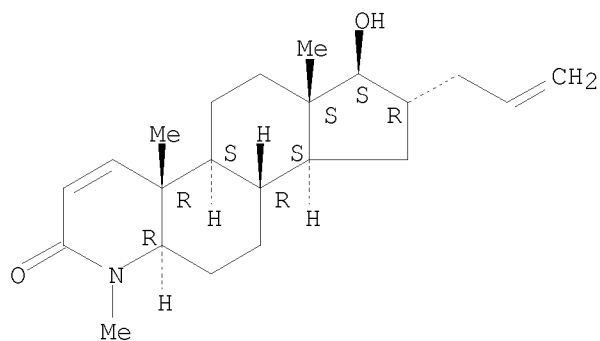
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(2-propen-1-yl)-,
 (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C22 H33 N O2

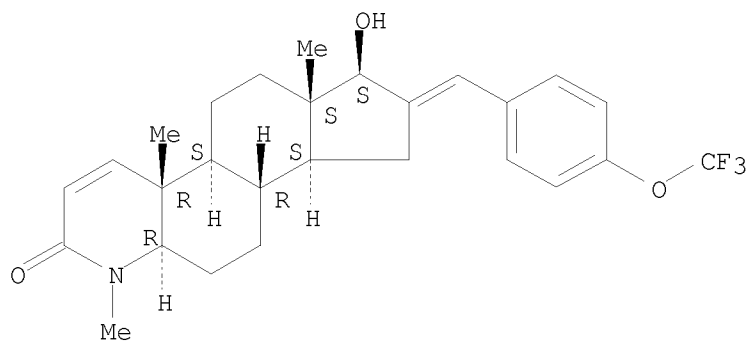
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[[4-
 (trifluoromethoxy)phenyl]methylene]-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C27 H32 F3 N O3

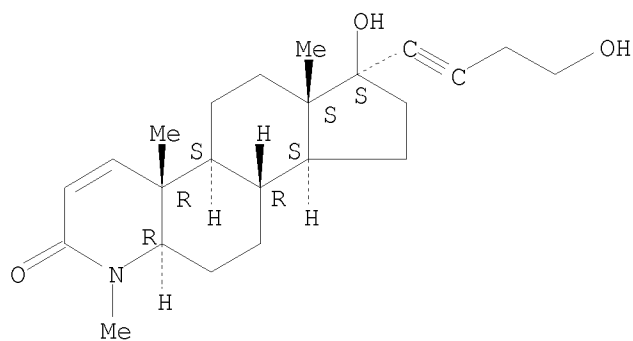
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-7-(4-hydroxy-1-butyn-1-yl)-1,4a,6a-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H33 N O3

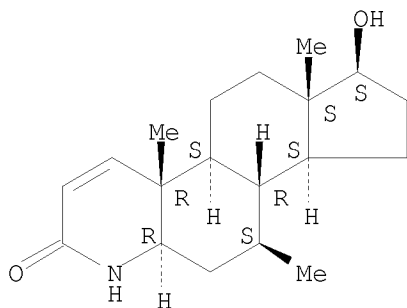
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-4a,6a,10-trimethyl-,
 (4aR,4bS,6aS,7S,9aS,9bR,10S,11aR)-
 MF C19 H29 N O2

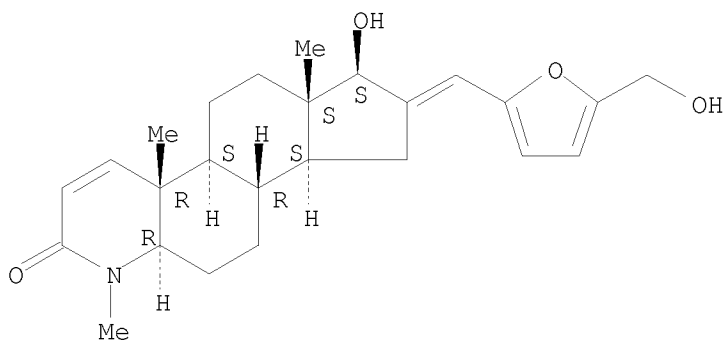
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-8-[[5-(hydroxymethyl)-2-furanyl]methylene]-
 1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C25 H33 N O4

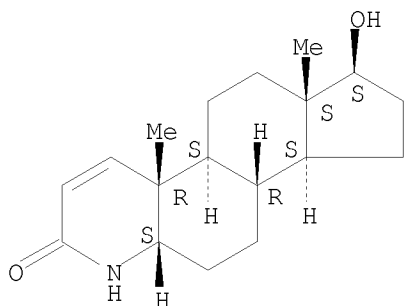
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-4a,6a-dimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aS)-
 MF C18 H27 N O2

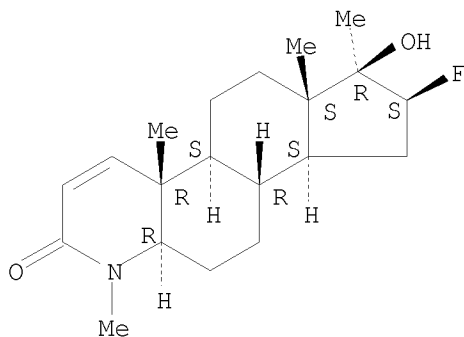
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-fluoro-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a,7-
 tetramethyl-, (4aR,4bS,6aS,7R,8S,9aS,9bR,11aR)-
 MF C20 H30 F N O2

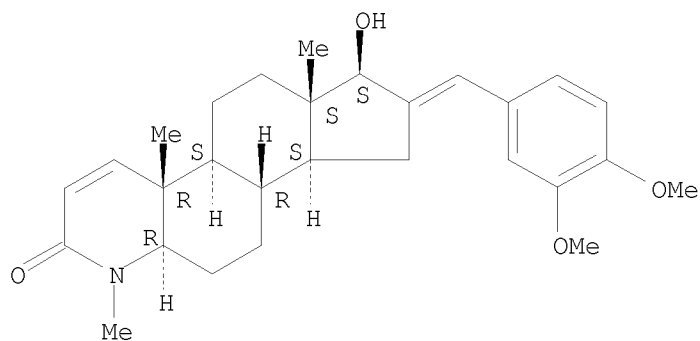
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(3,4-dimethoxyphenyl)methylene]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C28 H37 N O4

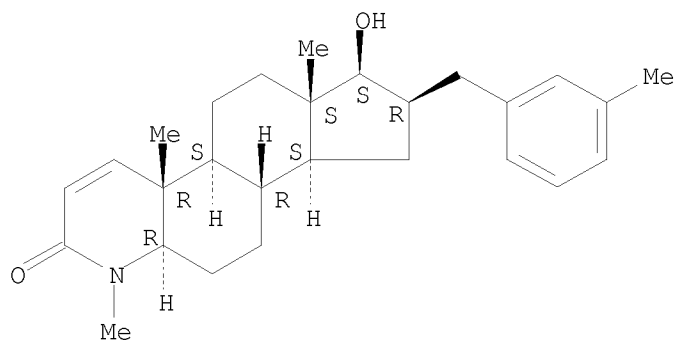
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-[(3-methylphenyl)methyl]-, (4aR,4bS,6aS,7S,8R,9aS,9bR,11aR)-
 MF C27 H37 N O2

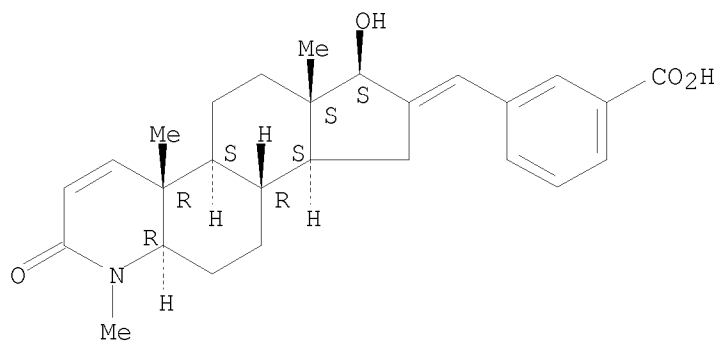
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN Benzoic acid, 3-[[[(4aR,4bS,6aS,7S,9aS,9bR,11aR)-1,2,4a,4b,5,6,6a,7,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-2-oxo-8H-indeno[5,4-f]quinolin-8-ylidene)methyl]-
 MF C27 H33 N O4

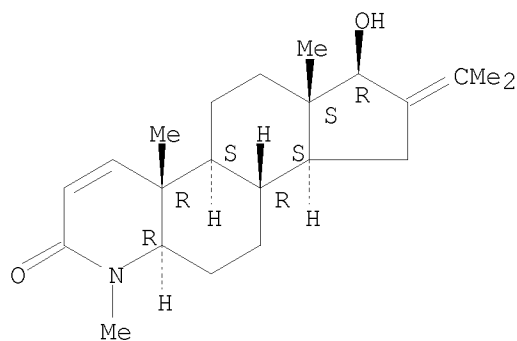
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(1-methylethylidene)-,
 (4aR,4bS,6aS,7R,9aS,9bR,11aR)-
 MF C22 H33 N O2

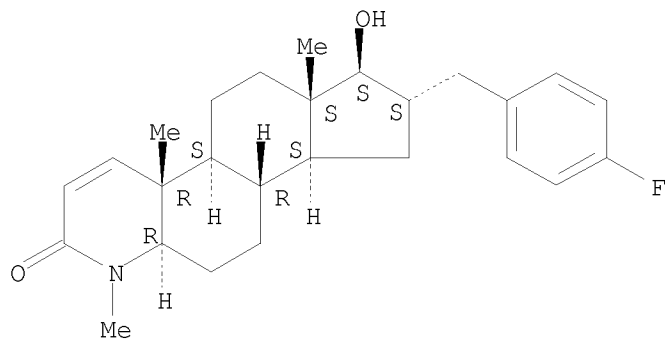
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-[(4-fluorophenyl)methyl]-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C26 H34 F N O2

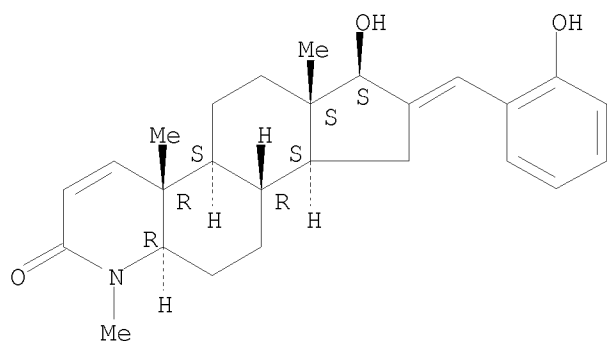
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-8-[(2-hydroxyphenyl)methylene]-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C26 H33 N O3

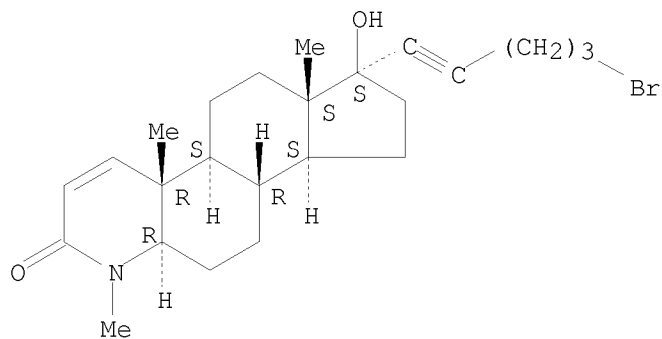
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 7-(5-bromo-1-pentyn-1-yl)-1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H34 Br N O2

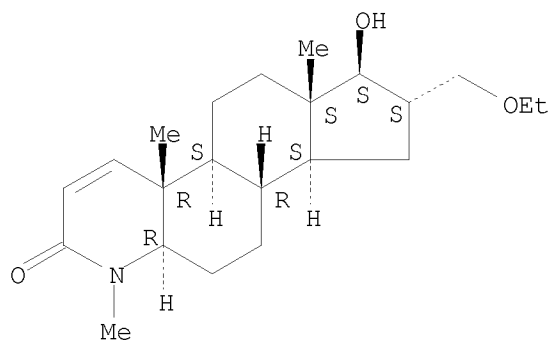
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 8-(ethoxymethyl)-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,8S,9aS,9bR,11aR)-
 MF C22 H35 N O3

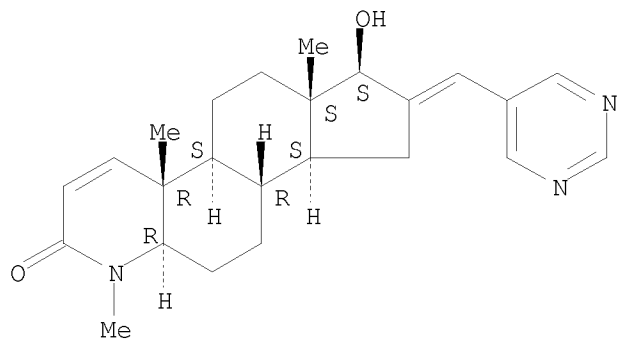
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-
 tetradecahydro-7-hydroxy-1,4a,6a-trimethyl-8-(5-pyrimidinylmethylene)-,
 (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C24 H31 N3 O2

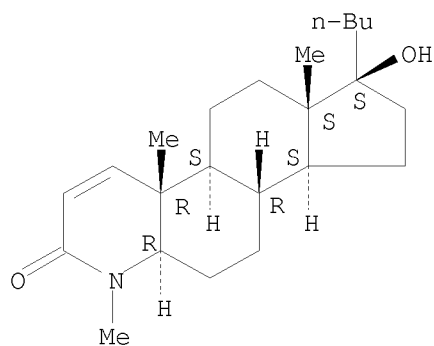
Absolute stereochemistry.
 Double bond geometry unknown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 103 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN
 IN 2H-Indeno[5,4-f]quinolin-2-one, 7-butyl-
 1,4a,4b,5,6,6a,7,8,9,9a,9b,10,11,11a-tetradecahydro-7-hydroxy-1,4a,6a-
 trimethyl-, (4aR,4bS,6aS,7S,9aS,9bR,11aR)-
 MF C23 H37 N O2

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=>
 Connection closed by remote host